

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017007.D
 Acq On : 08 Sep 2023 12:11
 Operator : MA/JU
 Sample : PB155068BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK068

Quant Time: Sep 09 00:43:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	372594	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1610472	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	985608	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2071988	20.000	ng/u1	-0.01
79) Chrysene-d12	21.498	240	1520030	20.000	ng/u1	-0.01
88) Perylene-d12	24.033	264	1447137	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	86011	8.732	ng/uL	0.00
4) Pyridine-d5	3.758	84	1004348	35.684	ng/u1	0.00
7) Phenol-d5	7.122	99	1180236	34.672	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	775424	36.418	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	880494	35.740	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	918006	35.399	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	440657	35.269	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	475801	35.130	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	827732	33.103	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1245794	34.533	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2710767	36.948	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3009760	35.547	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	449652	31.376	ng/u1	-0.01
60) Fluorene-d10	15.628	176	2297544	36.576	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	291002	23.338	ng/u1	-0.01
73) Anthracene-d10	17.504	188	3495849	36.739	ng/u1	0.00
81) Pyrene-d10	19.739	212	3769631	40.715	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.868	264	2900505	39.508	ng/u1	-0.01

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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